

Spontaneous Symmetry Breaking in Superfluid Helium-4

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We derive an analytical expression for a critical temperature of spontaneous symmetry breaking in a repulsive hard-core interacting Bose system. We show that the critical temperature of spontaneous symmetry breaking in a hard-core interacting Bose system is determined by the three physical parameters: the density of Bose liquid at absolute zero (ρ_0), the mass (m) and the hard sphere diameter (σ) of a boson. The formula that we have derived is $T_c = \rho_0 \pi \hbar^2 \sigma / m^2 k_B$. We report T_c of liquid helium-4 is 2.194 K, which is significantly close to the λ -temperature of 2.1768 K. The deviation between the predicted and experimental values of the λ -temperature is less than 1%.

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1. The year 2008 is the 100th anniversary of “Liquid Helium Year”, in which Heike Kamerlingh Onnes produced liquid helium. We are in a truly memorial moment in the history of low temperature physics. Liquid helium has been studied for the past 100 years, and there is no question that liquid helium is of central importance in this field. (I would like to turn readers’ attention to Ref. [1]. It presents the first complete English translation of the inaugural speech of Heike Kamerlingh Onnes at the University of Leiden in 1882. Although his speech is not related to helium, it is quite interesting and gives us a lesson. I believe that it attracts the interest of readers.)

In spite of great efforts, an important problem still remains to be solved. The problem is as follows. Liquid helium-4 undergoes a phase transition, known as the λ -transition, to a superfluid phase at the λ -temperature. The experimental value of the λ -temperature T_λ is approximately equal to 2.2 K. The problem is quite simple. Why is $T_\lambda \simeq 2.2$ K? What physical parameters determine the value of the λ -temperature? This is the “Why 2.2 K?” problem, which is the subject of the present paper. It is a long-standing dream in low temperature physics to derive a formula for the liquid helium-4 λ -temperature.

In 1938, Fritz London proposed that the λ -transition of liquid helium-4 probably has to be regarded as the Bose-Einstein condensation (BEC) [2]. He calculated a critical temperature of the BEC in an ideal Bose gas:

$$T_{BE} = \frac{2\pi\hbar^2}{mk_B} \left(\frac{n}{\zeta(3/2)} \right)^{2/3}, \quad (1)$$

where m is a boson mass, n is a number density of Bose gas and $\zeta(3/2) = 2.612375 \dots$. Thus, the Bose-Einstein temperature T_{BE} is determined by the two physical parameters, m and n . London reported T_{BE} of helium is 3.13 K [3], which is the same order of magnitude as the λ -temperature of 2.18 K. His proposal is quite important, because it was the first suggestion that the λ -transition was intrinsically related to the Bose-Einstein condensation. Recall that it was not clear in the 1930s whether the BEC was an actually physical phenomenon [3].

Although London’s work was revolutionary progress, there is a difference of approximately 0.95 K between his theoretical prediction and the experimental value. This deviation is approximately 44% and it is not negligible. It is widely believed that this discrepancy arises from the neglect of interactions between helium atoms in his theoretical calculation. This conjecture is quite reasonable because liquid helium is not an ideal Bose gas but a repulsive interacting Bose system. However, in general, it is difficult to consider interactions between atoms in analytical calculations of the BEC. Although this problem has been discussed [4], it remains very important.

In this paper, we study the “Why 2.2 K?” problem via an alternative approach. The key concept in our approach is *spontaneous symmetry breaking*. We derive an analytical formula for a critical temperature of spontaneous symmetry breaking in a repulsive hard-core interacting Bose system. We show that a critical temperature is determined by the three physical parameters: the density of Bose liquid at absolute zero (ρ_0), the mass (m) and the hard sphere diameter (σ) of a boson. Our formula predicts that a critical temperature of liquid helium-4 is 2.194 K, which is significantly close to the experimental value of the λ -temperature ($T_\lambda = 2.1768$ K [5]).

2. Spontaneous symmetry breaking is a symmetry breaking by the ground state of a system. The symmetry can be discrete or continuous, and be local or global. Spontaneous symmetry breaking is a phenomenon that occurs in many systems. In condensed matter physics, ferromagnetism is a primary example. For superconductivity, the importance of spontaneous symmetry breaking was firstly emphasized by Yoichiro Nambu [6]. In particle physics, the electroweak symmetry breaking is established in the Standard Model. In cosmology, a number of spontaneous symmetry breaking play an important role in the history of the early universe. Furthermore, the recently proposed *ghost condensation* [7] in the accelerating universe can be regarded as a kind of spontaneous symmetry breaking. Thus, spontaneous symmetry breaking is a very important concept in modern physics.

Spontaneous symmetry breaking is crucial in liquid helium-4. In short, *the λ -transition is the spontaneous symmetry breaking*. This is the basis of our approach. We derive a formula for a critical temperature of spontaneous symmetry breaking in the framework of effective field theory described by the order parameters. In the case of liquid helium-4, the order parameter is a macroscopic wavefunction φ , which is a one-component complex scalar field with both amplitude and phase. We begin with the theory described by this scalar field φ .

3. The Lagrangian density $\mathcal{L}(=\mathcal{K}-\mathcal{V})$ of the system is given by the nonrelativistic Goldstone model of the form

$$\mathcal{V} = -\mu\varphi^*\varphi + \frac{\lambda}{2}(\varphi^*\varphi)^2, \quad \lambda = \frac{2\pi\hbar^2\sigma}{m}. \quad (2)$$

The kinetic part $\mathcal{K} = i\hbar\varphi^*\partial_t\varphi + \hbar^2\varphi^*\nabla^2\varphi/2m$ is not important here, and therefore we concentrate on the potential \mathcal{V} . In Eq. (2), μ is the chemical potential, λ is the coupling constant for helium interactions and φ^* is the complex conjugate of φ . The interaction between helium atoms is repulsive at short distances (hard-core interactions). Therefore, the coupling constant $\lambda = 2\pi\hbar^2\sigma/m$ is positive, where m and σ are the mass and the hard sphere diameter of a helium atom, respectively.

We first consider the physical dimensions of parameters of the potential (2). In units of $\hbar = 1$ and $k_B = 1$ (\hbar is the reduced Planck's constant and k_B is the Boltzmann constant), for any physical quantity Q , its physical dimension $[Q]$ is written as the product of the dimensions of length and temperature:

$$[Q] = [\text{Length}]^\alpha[\text{Temperature}]^\beta, \quad (3)$$

where α and β are numerical constants. The physical dimension of the Lagrangian density is $[\mathcal{L}] = [\text{L}]^{-3}[\text{T}]$, and that of the scalar field φ is $[\varphi] = [\text{L}]^{-3/2}$. (Here $[\text{L}]$ and $[\text{T}]$ represent [Length] and [Temperature], respectively.) Hence it is straightforward to derive the physical dimensions of parameters in Eq. (2):

$$[\mu] = [\text{T}], \quad [\lambda] = [\text{L}]^3[\text{T}], \quad [m] = [\text{L}]^{-2}[\text{T}]^{-1}, \quad [\sigma] = [\text{L}]. \quad (4)$$

Note here that in the present case the coupling constant λ is a dimensionful parameter, in contrast to a relativistic case in which λ is dimensionless. These physical dimensions help us to understand the physics of liquid helium-4.

We next consider the symmetry of Lagrangian. The potential \mathcal{V} of Eq. (2) is invariant under the U(1) phase transformation of the field φ : $\varphi \rightarrow e^{i\theta}\varphi$, $\varphi^* \rightarrow e^{-i\theta}\varphi^*$. Furthermore, if this is a global transformation in which the transformation parameter θ does not depend on the space and time coordinates, the kinetic part \mathcal{K} is invariant under the U(1) phase transformation. Hence, the system described by the Lagrangian density \mathcal{L} has a global $\text{U}(1)\simeq\text{O}(2)$ symmetry. This global U(1) symmetry is spontaneously broken at low temperature as you see below.

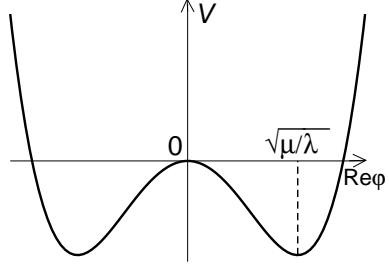


FIG. 1: The potential \mathcal{V} in the case that $\mu > 0$.

4. The ground state of a system is given by solving the condition $\partial\mathcal{V}/\partial\varphi = 0$. First, we consider the case that a chemical potential μ is *negative* ($\mu \leq 0$). In this case, there is only a trivial solution $\varphi = 0$. This solution is the stable ground state, because $\partial^2\mathcal{V}/\partial\varphi^*\partial\varphi|_{\varphi=0} \geq 0$. Hence the global U(1) symmetry is unbroken in this case. Second, we consider the case that μ is *positive* ($\mu > 0$). In this case, in contrast to the case that $\mu \leq 0$, there is a nontrivial solution in addition to a trivial solution $\varphi = 0$:

$$|\varphi| = \sqrt{\frac{\mu}{\lambda}}. \quad (5)$$

Fig. 1 shows the potential \mathcal{V} in the case that $\mu > 0$. In the present case, the trivial solution $\varphi = 0$ is unstable and it is not the ground state of the system. The nontrivial solution $|\varphi| = \sqrt{\mu/\lambda}$ is the stable ground state, because the curvature of the potential at two solutions satisfies

$$\left. \frac{\partial^2\mathcal{V}}{\partial\varphi^*\partial\varphi} \right|_{\varphi=0} < 0, \quad \left. \frac{\partial^2\mathcal{V}}{\partial\varphi^*\partial\varphi} \right|_{|\varphi|=\sqrt{\mu/\lambda}} > 0. \quad (6)$$

Therefore, if a chemical potential μ is *positive* ($\mu > 0$), the global U(1) symmetry is spontaneously broken. From above arguments, we reach a following view. A physical system undergoes a phase transition at a critical temperature T_c , if a chemical potential satisfies the conditions: $\mu \leq 0$ ($T \geq T_c$), $\mu > 0$ ($T < T_c$). Spontaneous symmetry breaking occurs only if a chemical potential becomes *positive*. Therefore, in contrast to the Bose-Einstein condensation, *spontaneous symmetry breaking never occurs in an ideal Bose gas*, in which a chemical potential is always *negative* ($\mu \leq 0$). This is a crucial difference between spontaneous symmetry breaking and the standard Bose-Einstein condensation.

5. Before deriving a formula for a critical temperature, we consider the temperature dependence of a chemical potential μ . Although the temperature dependence of μ is not necessary for deriving a formula, it is useful to understand a phase transition. For this reason, we consider the temperature dependence of μ in the following. The quantity $|\varphi|^2$ is equivalent to the number density n of superfluid. Therefore, the following relation is satisfied below the λ -temperature

$$|\varphi| = \sqrt{\frac{\mu}{\lambda}} = \sqrt{n} = \sqrt{\frac{\rho_s}{m}}, \quad (7)$$

where ρ_s is the superfluid density and m is a boson mass. From this relation, the superfluid density is given by

$$\rho_s = \frac{m\mu}{\lambda}. \quad (8)$$

This expression indicates that the temperature dependence of μ is obtained from that of ρ_s .

Here we take a minimal model:

$$\frac{\rho_s}{\rho} = 1 - \left(\frac{T}{T_\lambda} \right)^6, \quad (9)$$

where $\rho (= \rho_n + \rho_s)$ is the total density of liquid helium, and ρ_n, ρ_s are the normal fluid and superfluid densities, respectively. It should be emphasized that this model does not affect a formula for a critical temperature. Fig. 2 shows the superfluid density ratio as a function of temperature. Two functions $1 - (T/T_\lambda)^5$ and $1 - (T/T_\lambda)^7$ are plotted, for comparison. Fig. 2 shows that Eq. (9) and experimental values are consistent except near the λ -transition, at which the superfluid density ratio is of the form $(1 - T/T_\lambda)^{2/3}$. The form of Eq. (9) is the same as that of the Bose-Einstein condensation in an ideal Bose gas, $1 - (T/T_c)^{3/2}$. Although the exponent “6” has not yet been derived, we do not consider this problem here.

From Eqs. (8) and (9), we obtain the expression

$$\mu = \mu_0 \left(1 - \left(\frac{T}{T_\lambda} \right)^6 \right), \quad \mu_0 = \frac{\lambda \rho_0}{m}, \quad (10)$$

where μ_0 is a chemical potential at absolute zero and ρ_0 is the total density at absolute zero. We have used the approximation $\rho \simeq \rho_0$ in Eq. (10), because the total density ρ is approximately a constant below the λ -temperature. Therefore, Eq. (10) is a good approximation at low temperature ($T < 1.7$ K). Fig. 3 shows the schematic temperature dependence of the potential \mathcal{V} . For $T > T_\lambda$, the potential has only one minimum at $\varphi = 0$ and the curvature at $\varphi = 0$ is positive. When $T = T_\lambda$, the curvature at the minimum is zero, $\partial^2 \mathcal{V} / \partial \varphi^* \partial \varphi|_{\varphi=0} = 0$. For $T < T_\lambda$, the curvature at $\varphi = 0$ is negative and the scalar field rolls down to the $|\varphi| \neq 0$ minimum. Thus, there is no barrier of potential between the minimum at $\varphi = 0$ and $|\varphi| \neq 0$, indicating a second-order phase transition.

6. We now derive a formula for a critical temperature of spontaneous symmetry breaking. When $T = 0$, the field φ is located at the nonzero minimum $|\varphi| \neq 0$. Hence the global U(1) symmetry is broken in this phase. As the temperature increases above a critical temperature, the potential has only one minimum at $\varphi = 0$ and the global U(1) symmetry is recovered. From these observations, we reach a following conclusion; *the depth of the potential well at absolute zero determines a critical temperature*. When $T = 0$, the depth of the potential well is given by

$$-\mathcal{V}(|\varphi| = \sqrt{\mu_0/\lambda}) = \frac{\mu_0 n_0}{2}, \quad (11)$$

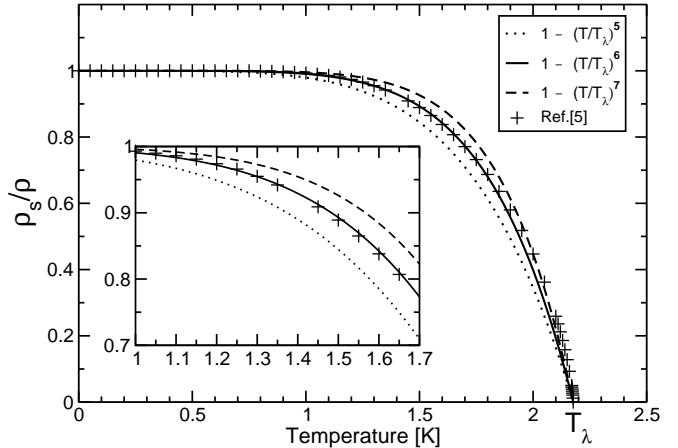


FIG. 2: Superfluid density ratio for liquid helium-4 as a function of temperature. The inset shows the temperature region $1.0 \text{ K} \leq T \leq 1.7 \text{ K}$. Key: (dotted line) $1 - (T/T_\lambda)^5$; (solid line) $1 - (T/T_\lambda)^6$; (dashed line) $1 - (T/T_\lambda)^7$; (+) the data from Ref. [5].

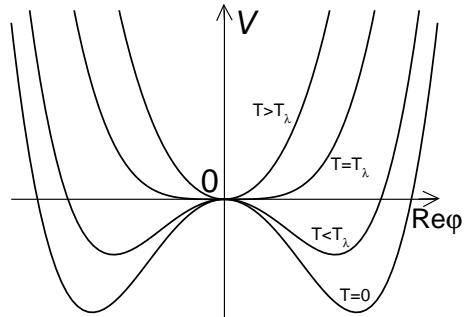


FIG. 3: The schematic temperature dependence of the potential \mathcal{V} for a second-order phase transition.

where $n_0 = \rho_0/m$ is the number density of Bose liquid at absolute zero. This quantity $\mu_0 n_0/2$ represents the required energy density to recover the global U(1) symmetry of the system. Therefore, the corresponding thermal energy density $k_B T_c n_0$ is equivalent to $\mu_0 n_0/2$. Consequently, we obtain the following relation

$$k_B T_c = \frac{\mu_0}{2}. \quad (12)$$

Fig. 4 shows the present situation. From above arguments, we have the following relations:

$$k_B T_c = \frac{\mu_0}{2}, \quad \frac{\rho_0}{m} = \frac{\mu_0}{\lambda}, \quad \lambda = \frac{2\pi\hbar^2\sigma}{m}. \quad (13)$$

Therefore, a critical temperature of spontaneous symmetry breaking in a repulsive interacting Bose system is:

$$T_c = \frac{\rho_0 \pi \hbar^2 \sigma}{m^2 k_B}, \quad (14)$$

where ρ_0 is the density of Bose liquid at absolute zero, m and σ are the mass and the hard sphere diameter of a boson, respectively. Substituting the values of

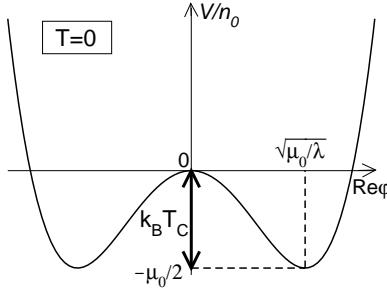


FIG. 4: The absolute zero potential per number density.

liquid helium-4 into Eq. (14), $\rho_0 = 0.1451 \text{ g/cm}^3$ [5], $m = 6.6465 \times 10^{-24} \text{ g}$, and $\sigma = 2.639 \text{ \AA}$ [8, 9], we obtain the value $T_c = 2.194 \text{ K}$. This prediction is significantly close to the experimental value of the λ -temperature 2.1768 K [5]. The deviation Δ between the predicted and experimental values of the λ -temperature is less than 1%:

$$\Delta \equiv 100 \times \frac{T_c^{(\text{theory})} - T_\lambda^{(\text{exp})}}{T_\lambda^{(\text{exp})}} \simeq 0.8\% < 1\%. \quad (15)$$

Thus, our prediction is much closer to the experimental value than London's prediction. It is because in the present approach the repulsive interactions between helium atoms are included in the derivation of a formula. While the Bose-Einstein temperature T_{BE} in an ideal Bose gas is determined by the *two* physical parameters, our formula for a critical temperature, Eq. (14), is determined by the *three* physical parameters: the density of superfluid (ρ_0), the mass (m) and the hard sphere diameter (σ) of a boson. The additional physical parameter σ includes information about interactions, and it has significantly improved the theoretical prediction for a critical temperature. The results are summarized in Table I.

Finally, we comment on the hard sphere diameter σ . Although ρ_0 and m have been determined precisely, the uncertainty of σ is relatively large. The hard sphere diameter σ represents the point at which the interatomic potential $V(r)$ is zero ($V(\sigma) = 0$), and the distance r_m represents the point at which the potential $V(r)$ is minimum ($\partial V(r)/\partial r|_{r=r_m} = 0$). Although the value of r_m has been reported by many groups [10], that of σ has been less reported. Furthermore, the relation $r_m = 2^{1/6}\sigma$ of the (12,6) Lennard-Jones potential model is not appli-

TABLE I: A comparison between BEC and SSB. Right row is given in this work. The deviation Δ is defined by Eq. (15).

System	ideal Bose gas	interacting Bose liquid
Key concept	BEC	SSB
T_c	$2\pi\hbar^2(n/\zeta(3/2))^{2/3}/mk_B$	$\rho_0\pi\hbar^2\sigma/m^2k_B$
parameter#	2	3
T_c (${}^4\text{He}$)	3.13 K	2.194 K
Δ	44 %	0.8 %

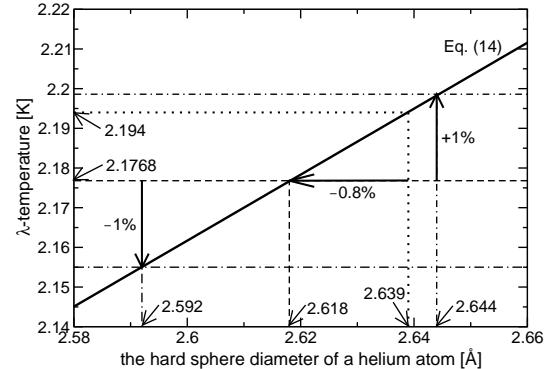


FIG. 5: The λ -temperature as a function of the hard sphere diameter σ of a helium atom.

cable to the present study, because it is not sufficiently a good approximation. For these reasons, we plot the λ -temperature as a function of σ . Fig. 5 shows that the deviation between the predicted and experimental values is less than $\pm 1\%$ for the region $2.592 \text{ \AA} \leq \sigma \leq 2.644 \text{ \AA}$.

7. In conclusion, we have derived an analytical formula for a critical temperature of spontaneous symmetry breaking in a repulsive interacting Bose system. The formula that we have derived, Eq. (14), is a simple analytical expression to predict $T_\lambda \simeq 2.2 \text{ K}$. We hope that this work contributes to progress of low temperature physics.

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